

Ground-state phase diagram of a half-filled one-dimensional extended Hubbard model

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The density-matrix renormalization group is used to study the phase diagram of the one-dimensional half-filled Hubbard model with on-site (U) and nearest-neighbor (V) repulsion, and hopping t . A critical line $V_c(U) \approx U/2$ separates a Mott insulating phase from a charge-density-wave phase. The formation of bound charge excitations for $V > 2t$ changes the phase transition from continuous to first order at a tricritical point $U_t \approx 3.7t, V_t = 2t$. A frustrating effective antiferromagnetic spin coupling induces a bond-order-wave phase on the critical line $V_c(U)$ for $U_t < U \lesssim 7t$.

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The properties of quasi-one-dimensional materials have been extensively studied in recent years [1, 2, 3]. These materials exhibit rich phase diagrams and display unusual optical properties due to the combination of reduced dimensionality and strong electronic correlations. Consequently, much effort has been devoted to understanding the ground-state and optical properties of theoretical one-dimensional correlated electron systems such as the half-filled Hubbard model with on-site (U) and nearest-neighbor (V) repulsion and hopping term t . Nevertheless, the ground state phase diagram of this model is still controversial [4, 5, 6, 7, 8, 9, 10]. It is known [4] that the system is a Mott insulator for $U \gtrsim 2V$ and a charge density wave (CDW) insulator for $U \lesssim 2V$. The quantum phase transition is continuous at weak coupling ($U, V \ll t$) and first order at strong coupling ($U, V \gg t$). Numerically [4, 5, 10], one finds that the order of the transition changes at a tricritical point (U_t, V_t) with $V_t/t \approx 1.5 - 2.5$, but this feature is not well understood. Recently, it has been proposed [8, 9, 10] that a bond-order-wave (BOW) phase exists between the Mott and CDW phases up to the tricritical point. Concurrently, the optical properties of this model have been determined in the Mott insulating phase [11, 12, 13, 14, 15]. In particular, it has been found that the lowest optical excitations consist of a pair of independent charge excitations for $V \leq 2t$, while they are bound states for $V > 2t$. Surprisingly, the remarkable proximity of the tricritical point to the boundary between bound and free charge excitations has not been noticed until now.

Here I investigate the ground-state phase diagram of the half-filled one-dimensional extended Hubbard model in the repulsive regime using the density-matrix renormalization group (DMRG) [16]. I show that the nature of the low-lying charge excitations determines the order of the transition and the position of the tricritical point. A BOW phase is found only at intermediate coupling on the critical line $V_c(U)$ between Mott and CDW phases.

The model is defined by the Hamiltonian

$$\hat{H} = -t \sum_{l,\sigma} \left(\hat{c}_{l,\sigma}^\dagger \hat{c}_{l+1,\sigma} + \hat{c}_{l+1,\sigma}^\dagger \hat{c}_{l,\sigma} \right)$$

$$+ U \sum_l \left(\hat{n}_{l,\uparrow} - \frac{1}{2} \right) \left(\hat{n}_{l,\downarrow} - \frac{1}{2} \right) + V \sum_l (\hat{n}_l - 1)(\hat{n}_{l+1} - 1). \quad (1)$$

Here $\hat{c}_{l,\sigma}^\dagger, \hat{c}_{l,\sigma}$ are creation and annihilation operators for electrons with spin $\sigma = \uparrow, \downarrow$ at site $l = 1, \dots, N$, $\hat{n}_{l,\sigma} = \hat{c}_{l,\sigma}^\dagger \hat{c}_{l,\sigma}$, and $\hat{n}_l = \hat{n}_{l,\uparrow} + \hat{n}_{l,\downarrow}$. I exclusively consider systems with an even number N of sites. At half filling the number of electrons equals N . The interaction is repulsive $U \geq V \geq 0$. The Hamiltonian (1) has a particle-hole symmetry and a spatial reflection symmetry. Thus, each eigenstate has a well-defined parity under charge conjugation ($P_c = \pm 1$) and belongs to one of the two irreducible representations, A_g or B_u , of the reflection symmetry group. The ground state belongs to the symmetry subspace $A_g^+ \equiv (A_g, P_c)$ and optically excited states belong to the symmetry subspace $B_u^- \equiv (B_u, -P_c)$ because the current operator is antisymmetric under charge conjugation and spatial reflection [14].

DMRG [16] is known to be a very accurate numerical method for one-dimensional quantum systems with short-range interactions such as the Hamiltonian (1). Here I use the finite-system DMRG algorithm to calculate eigenenergies and static properties at low-energy. Spin gaps $E_s = E_0(N, 1) - E_0(N, 0)$ and single-particle charge gaps $E_c = 2[E_0(N+1, 1/2) - E_0(N, 0)]$ are derived from the ground state energies $E_0(N_e, S_z)$ for a given number of electron N_e and a given spin S_z . I also use a symmetrized DMRG [15, 17] method to calculate the energy and static properties of the lowest eigenstates in the B_u^- symmetry subspace.

All DMRG methods have a truncation error which is reduced by increasing the number m of density-matrix eigenstates kept in the renormalization procedure [16]. To achieve a greater accuracy and to obtain error estimates I extrapolate DMRG results to the limit of vanishing discarded weight P_m . For eigenenergies DMRG errors vanish linearly with P_m , while for other quantities DMRG errors usually scale as $(P_m)^\alpha$ with $0 < \alpha \leq 1$. The largest number of density-matrix eigenstates used in this work is $m = 1200$ and truncation errors are negli-

TABLE I: Critical nearest-neighbor repulsion $V_c(U)$, single-particle charge gap E_c at (U, V_c) , and transition type for several values of U . Numbers in parenthesis are estimated errors.

U/t	$V_c(U)/t$	E_c/t	transition order
2	1.125 (25)	0	continuous
3	1.640 (10)	0	continuous
4	2.150 (10)	0.05	first order
5	2.665 (5)	0.11	first order
6	3.155 (5)	0.15	first order
8	4.141 (4)	0.45	first order
12	6.115 (5)	2.0	first order
40	20.041 (4)	29.3	first order

gible for all results presented here. DMRG calculations have always been carried out for several sizes N in order to check finite-size effects and, if necessary, results have been extrapolated to the thermodynamic limit $N \rightarrow \infty$. The largest system size used in this work is $N = 1024$.

For $V \lesssim U/2$ the model (1) describes a Mott insulator. [This phase is often called a spin-density-wave (SDW) phase because antiferromagnetic spin correlations decay algebraically in the ground state.] There is no broken symmetry and the ground state is nondegenerate. The charge gap E_c is finite but the spin gap E_s vanishes in the thermodynamic limit. In this Mott insulating phase the lowest eigenstate in the B_u^- symmetry sector always contributes to the optical spectrum. Therefore, the difference between the B_u^- and A_g^+ ground-state energies corresponds to the optical gap E_{opt} . For $V \gtrsim U/2$ the system is in a long-range ordered CDW phase with a doubly degenerate ground state in the thermodynamic limit. In this phase both charge and spin gaps are finite. The CDW order parameter $0 < |m_e| \leq 1$ gives the amplitude of the ground-state density modulation $\langle \hat{n}_l \rangle = 1 + (-1)^l m_e$. Such a CDW breaks the charge-conjugation symmetry and the spatial reflection symmetry. Thus one of the two degenerate ground states belongs to the A_g^+ symmetry subspace and the other one to the B_u^- subspace. Obviously, this consideration shows that the low-energy B_u^- excitations in the Mott phase play an important role in the Mott-CDW transition.

The transition from the CDW phase to the Mott phase can be investigated with DMRG as done previously for the CDW-metal transition in the Holstein model [18]. More precisely, I have investigated the lowest eigenstates to determine the ground-state degeneracy and symmetry, and I have calculated the electronic staggered susceptibility to check if the ground state has long-range CDW order. Some results for the critical line $V_c(U)$ determined with this approach are given in Table I. These results agree quantitatively with recent quantum Monte Carlo (QMC) simulations [10] and the Gaussian transition line in Ref. 8, although the interpretation of this phase boundary is completely different (see below). These values are also in good agreement with strong-

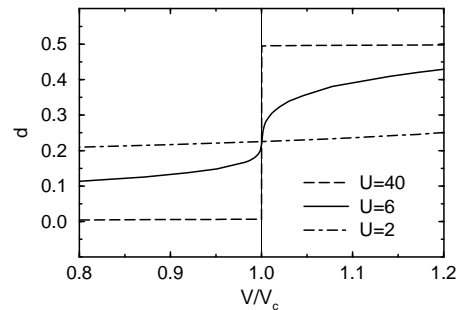


FIG. 1: Ground-state double occupancy d versus V close to the critical coupling $V_c(U)$ for several U (in units of t).

coupling perturbation theory [6] down to $U = 6t$ and the results of early numerical simulations [4, 5]. This confirms the accuracy of the present approach and I will not elaborate further on the location of this boundary.

To determine the order of the transition on the critical line $V_c(U)$, one can examine the derivatives of the ground-state energy per site with respect to the interaction parameters U and V . Using the Hellmann-Feynman theorem one shows that the derivatives are given (up to known constants) by ground-state expectation values of the double occupancy and nearest-neighbor density-density operators, respectively. These expectation values can be calculated accurately with DMRG. Figure 1 shows the average ground-state double occupancy

$$d = \frac{1}{N} \sum_l \langle \hat{n}_{l\uparrow} \hat{n}_{l\downarrow} \rangle \quad (2)$$

versus V close to the critical coupling $V_c(U)$. For $U = 40t$, d jumps from almost zero to slightly less than 0.5 at V_c and thus the transition is clearly first order, while it is continuous for $U = 2t$. For $U = 6t$ the derivative of d with respect to V evolves into a δ -function at $V = V_c$ as the system size N increases. Therefore, d is discontinuous at $V_c(U)$ and the transition is also first order for $U = 6t$. With this method I have found that the transition at $V_c(U)$ is first order from the strong-coupling limit ($U, V \gg t$) down to at least $U = 4t$ ($V_c \approx 2.15t$) and is continuous from the weak-coupling limit ($U, V \ll t$) up to at least $U = 3.5t$ ($V_c \approx 1.9t$).

An investigation of the low-lying charge excitations in the Mott insulating phase allows us to understand the existence of a tricritical point and to determine its position. Figure 2 shows the evolution of the charge gap E_c and the optical gap E_{opt} for three different couplings U . I have always found that the charge gap E_c varies continuously and goes through a minimum at $V_c(U)$ as V is increased for fixed U . For weak coupling ($U \leq 3t$) the gap vanishes smoothly on the critical line. For stronger coupling ($U \geq 4$) the slope of E_c with respect to V becomes very large and is discontinuous at V_c . In this regime the charge gap remains finite at the transition (see Table I).

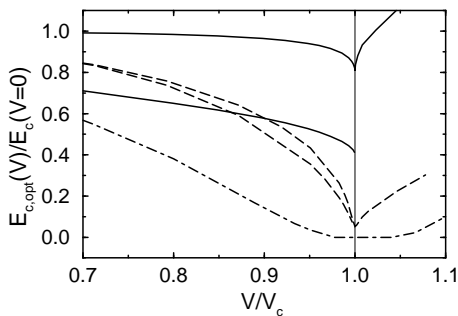


FIG. 2: Charge gap E_c (upper line) and optical gap E_{opt} (lower line) versus V close to the critical coupling $V_c(U)$ for $U/t = 40$ (solid), 6 (dashed), and 2 (dot-dashed). For $U = 2t$, $E_{\text{opt}} = E_c$. E_{opt} is not shown in the CDW phase ($V > V_c$).

In a Mott insulator, elementary charge excitations can be understood as spinless bosons in Hubbard bands. Excited states in the B_u^- subspace always consist of an even number of such elementary excitations with zero total charge [11, 13, 14]. DMRG calculations combined with analytical methods [14, 15] reveal their properties. For $V \leq 2t$ the low-energy B_u^- excited states consist of two independent elementary charge excitations and the optical gap equals the charge gap in the thermodynamic limit. For $V > 2t$, however, the lowest B_u^- excitations are bound states starting at an energy $E_{\text{opt}} < E_c$ (see Fig. 2). Just above $2t$ they are excitons made of two elementary charge excitations. Close to the critical line $V_c(U) \approx U/2$ they become finite-size “droplets” of the CDW phase, (i.e., a bound state of several elementary charge excitations). The excitation energy E_{opt} of the lowest CDW droplet remains finite at the transition $V_c(U)$ at least for $U \geq 5$ (see also Fig. 2). CDW droplet sizes increase sharply as V approaches the critical coupling $V_c(U)$ for fixed U and reach a finite value ξ_c at the critical line. This critical droplet size ξ_c diverges for $U \rightarrow \infty$ but tends to 2 (corresponding to an exciton) if V_c approaches $2t$.

Combining all results, one discerns two distinct regimes in the Mott phase as the critical line $V_c(U)$ is approached. If $V_c \leq 2t$, the transition is continuous, the B_u^- excitations are made of free charge excitations and become gapless at V_c . If $V_c > 2t$, the transition is first order, the low-lying B_u^- excitations are bound states and remain gapped at V_c . Therefore, I believe that the tricritical point is precisely located at the intersection of the $V = 2t$ boundary with the quantum critical line $V_c(U) \approx U/2$, i.e., at $U_t \approx 3.7t$, $V_t = 2t$.

Adapting Hirsch’s analysis [4], one can now qualitatively explain why the transition from the Mott to the CDW phase changes from first order to continuous. If $V_c(U) > 2t$, CDW droplets with sizes $\xi \geq \xi_c$ are energetically favored when V becomes larger than $V_c(U)$, and the system tunnels to the CDW phase by nucleation. On the critical line $V_c(U)$ the creation energy E_{opt} and crit-

ical size ξ_c of the lowest CDW droplet diminishes with U . At $V_c = 2t$, CDW droplets (with sizes $\xi \rightarrow 2$) become unstable and split into independent gapless elementary charge excitations. For $V_c(U) \leq 2t$, it is then advantageous to create many such excitations as V is increased above $V_c(U)$ and the transition becomes continuous.

In the field theory approach to the lattice model (1) the limit between free and bound charge excitations corresponds to a parameter $K_\rho = 1/2$ for a half-filled system [13, 14]. Field-theoretical predictions for the low-energy optical spectrum agree perfectly with dynamical DMRG calculations [14, 15]. Thus the tricritical point probably corresponds to $K_\rho = 1/2$. In Ref. 10, it has also been found that K_ρ is close to $1/2$ in this region.

Recently, Nakamura [8] has proposed that a narrow region with long-range BOW order exists between the CDW phase and the Mott phase at weak to intermediate coupling. The BOW region vanishes at the tricritical point, beyond which there is a direct first-order transition from the Mott phase to the CDW phase. In this scenario the critical line $V_c(U)$ calculated above would correspond to the BOW-CDW phase boundary for $U < U_t$. The existence of long-range BOW order in the critical region has been confirmed by QMC simulations [10].

In a BOW phase the ground state is two-fold degenerate and both charge and spin gaps are finite. A long-range ordered BOW corresponds to a finite staggered bond order

$$\frac{1}{2} \sum_{\sigma} \langle \hat{c}_{l,\sigma}^+ \hat{c}_{l+1,\sigma} + \hat{c}_{l+1,\sigma}^+ \hat{c}_{l,\sigma} \rangle = p_0 + (-1)^l \delta \quad (3)$$

with $\delta \neq 0$ in the thermodynamic limit. To distinguish the BOW phase from the Mott phase I have calculated the spin gap E_s and the bond order parameter δ using DMRG. After extrapolation to vanishing truncation errors P_m and to the thermodynamic limit $N \rightarrow \infty$, this approach gives spin gaps with an accuracy of $10^{-3}t$ or better and allows one to detect a bond order alternation as small as $|\delta| = 0.01$. (It appears numerically that δ either vanishes as $1/\sqrt{N}$ or tends to a finite value with finite-size corrections scaling as $1/N$.) I have found a BOW ground state only in a very narrow region adjacent to the critical line $V_c(U)$ for intermediate coupling $4t \leq U \leq 6t$. For $U = 4t$ and $V = 2.14t$, I have obtained $\delta = 0.08$ in quantitative agreement with QMC simulations [10]. However, the BOW phase extends to significantly stronger coupling than reported in Ref 10. For instance, there is a BOW ground state with $\delta \approx 0.12$ for $U = 6t$ and $V = 3.145t$. I have not observed any BOW ground state at weak coupling ($U \leq 3t$), at strong coupling ($U \geq 8t$), and for $V \leq U/2$. Actually, as BOW ground states are found only within $0.02t$ of the critical line $V_c(U)$, I believe that the BOW phase exists only on this critical line for intermediate coupling U starting from the tricritical point $U_t \approx 3.7t$ up to an upper limit which is smaller than $8t$. Field theory [7] also suggests

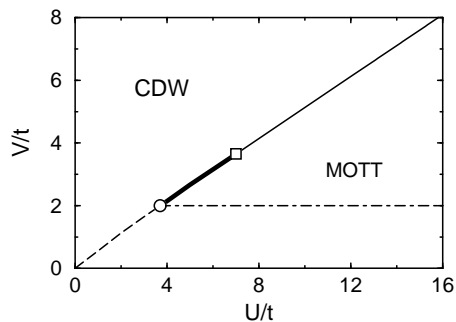


FIG. 3: Schematic ground-state phase diagram: the transition is continuous at weak coupling (dashed line) and first order at strong coupling (both solid lines). A circle marks the tricritical point. The dot-dashed line is the boundary between free and bound charge excitations in the Mott phase. The thick solid line indicates the BOW phase and a square marks the transition from a “spin-liquid” to a dimerized spin ground state on the critical line.

that a BOW phase can occur only on the boundary between CDW and Mott phases in the model (1). Even in the Hartree-Fock approximation, one finds a BOW phase only on the critical line $U = 2V$ between a SDW phase and a CDW phase [19].

The occurrence of the BOW phase on the critical line for $U \lesssim 8t$ can be understood as the result of increasing frustration in the spin degrees of freedom. On this critical line, low-energy charge excitations are dispersionless and thus ineffective in a degenerate CDW-Mott ground state. A strong-coupling expansion up to 4th order in t/U shows that the spin properties are determined by an effective Heisenberg Hamiltonian with nearest- (J_1) and next-nearest-neighbor (J_2) antiferromagnetic exchange couplings [6]. This strong-coupling perturbation analysis gives accurate results for the critical line $V_c(U)$ (and thus is expected to be valid) down to $U = 6t$. The ratio J_2/J_1 is strongly enhanced by the nearest-neighbor repulsion V . It is known [20, 21] that the frustration due to the J_2 coupling causes a quantum phase transition from a “spin-liquid” ground state for $4J_2 \lesssim J_1$ to a dimerized spin ground state for $4J_2 \gtrsim J_1$ in the Heisenberg model. On the critical line $V_c(U)$ of the extended Hubbard model, this corresponds to the appearance of a BOW phase (driven by the spin dimerization) for $U \lesssim 7t$, in agreement with my DMRG results, if one uses the values of J_1 and J_2 given by the 4th-order perturbation expansion [6]. Away from the critical line [$V < V_c(U)$], one observes only algebraically decreasing BOW correlations because charge fluctuations, which are no longer dispersionless, destroy the local spin moments and thus prevent any long-range spin order.

Figure 3 shows the phase diagram found with DMRG. This phase diagram disproves recent speculations about an extended BOW phase from weak coupling up to the tricritical point [8, 9, 10] but is compatible with the re-

sults of most previous investigations of this problem (see Refs. 4 to 7 and references therein).

In summary, I have studied the ground-state phase diagram of the one-dimensional half-filled extended Hubbard model for $U, V > 0$ using DMRG. I have shown that the phase diagram is determined by 3 mechanisms: First, the competition between the on-site (U) and nearest-neighbor (V) repulsion is responsible for a quantum critical transition from a Mott to a CDW ground state at $V \approx U/2$. Second, the competition between the hopping term t and the nearest-neighbor repulsion V is responsible for the transformation of free charge excitations into bound states (CDW droplets) at $V = 2t$ in the Mott phase and thus changes the Mott-CDW transition from continuous to first order. Third, a frustrating effective antiferromagnetic spin coupling leads to a dimerized spin ground state (and thus to a BOW phase) on the Mott-CDW critical line at intermediate coupling.

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